In the Claims:

1. (Original) A compound of formula (I):

$$Ar^{2}-Y-Ar^{1}$$

$$R^{3}$$

$$R^{2}$$

$$R^{3}$$

$$R^{1}$$

$$R^{1}$$

$$R^{1}$$

$$R^{2}$$

wherein

A and B represent the groups $-(CH_2)_m$ - and $-(CH_2)_n$ - respectively; R^1 represents C_{1-6} alkyl;

 R^2 represents hydrogen, halogen, hydroxy, cyano, nitro, hydroxyC₁₋₆alkyl, trifluoromethyl, trifluoromethoxy, C₁₋₆alkyl, C₁₋₆alkoxy, -(CH₂)_pC₃₋₆cycloalkyl, -(CH₂)_pOC₃₋₆cycloalkyl, -COC₁₋₆alkyl, -SO₂C₁₋₆alkyl, -SOC₁₋₆alkyl, -S-C₁₋₆alkyl, -CO₂C₁₋₆alkyl, -CO₂NR⁴R⁵, -(CH₂)_pNR⁴R⁵, -(CH₂)_pNR⁴COR⁵, an optionally substituted aryl group, an optionally substituted heterocyclyl group;

R³ represents hydrogen or C₁₋₆alkyl;

Ar¹ represents an optionally substituted heteroaryl group;

Ar² represents an optionally substituted phenyl or an optionally substituted heteroaryl group;

Y represents a bond, -O-, -C₁₋₆alkyl-, -CR⁶R⁷X-, -XCR⁶R⁷-, -NR⁸CO- or -CONR⁸-; X represents oxygen, sulfur, -SO- or -SO₂-;

R⁴ and R⁵ each independently represent hydrogen or C₁₋₆alkyl or, together with the nitrogen or other atoms to which they are attached, form an azacycloalkyl ring or an oxo-substituted azacycloalkyl ring;

 R^6 and R^7 each independently represent hydrogen, $C_{1\text{-}6}$ alkyl or fluoro;

R⁸ represents hydrogen or C₁₋₆alkyl;

m and n independently represent an integer selected from 1 and 2; p independently represents an integer selected from 0, 1, 2 and 3; or a pharmaceutically acceptable salt, solvate or pharmaceutically acceptable derivative thereof.

- 2. (Original) A compound of formula (I) which is
- 5-(4-Chlorophenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(3-Methoxyphenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(4-Methoxyphenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(3,4-Difluorophenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(2,4-Difluorophenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(3-Chlorophenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(3-Fluorophenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(4-Trifluoromethylphenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(3-Trifluoromethylphenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(4-Fluorophenyl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(4-Fluorophenyl)-thiophene-2-sulfonic acid (3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(4-Chlorophenyl)-thiophene-2-sulfonic acid (3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)amide;
- 5-(4-Chloro-2-methylphenyl)-thiophene-2-sulfonic acid (2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)-amide;
- 5-Isoxazol-3-yl-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)-amide;
- 5-(2-Methylthiazol-5-yl)-thiophene-2-sulfonic acid (8-methoxy-3-methyl-2,3,4,5-tetrahydro-1*H*-benzo[*d*]azepin-7-yl)-amide;

- [2,3']Bithiophenyl-5-sulfonic acid (2,3,4,5-tetrahydro-1H-benzo[d]azepin-7-yl)-amide;
- [2,3']Bithiophenyl-5-sulfonic acid (3-methyl-2,3,4,5-tetrahydro-1H-benzo[d]azepin-7-yl)-amide;
- 5-(4-Chlorophenyl)thiophene-2-sulfonic acid (8-dimethylamino-3-methyl-2,3,4,5-tetrahydro-1*H*-3-benzazepin-7-yl)amide;
- 5-(4-Fluorophenyl)thiophene-2-sulfonic acid (8-dimethylamino-3-methyl-2,3,4,5-tetrahydro-1*H*-3-benzazepin-7-yl)amide;
- 5-(2,4-Difluorophenyl)thiophene-2-sulfonic acid (8-dimethylamino-3-methyl-2,3,4,5-tetrahydro-1*H*-3-benzazepin-7-yl)amide; and
- 5-(3,4-Difluorophenyl)thiophene-2-sulfonic acid (8-dimethylamino-3-methyl-2,3,4,5-tetrahydro-*1H*-3-benzazepin-7-yl)amide.
- 3. (Currently Amended) A pharmaceutical composition comprising a compound of formula (I) as claimed in claim 1 or claim 2 or a pharmaceutically acceptable derivative thereof and a pharmaceutically acceptable carrier therefor.

Claims 4-8. (Cancelled)

- 9. (Currently Amended) A method of treating a condition which requires modulation of dopamine receptors which comprises administering to a mammal in need thereof an effective amount of a compound of formula (I) according to claim 1-or elaim 2.
- 10. (Original) A method of treating a condition according to claim 9 wherein the condition is schizophrenia or substance abuse.